

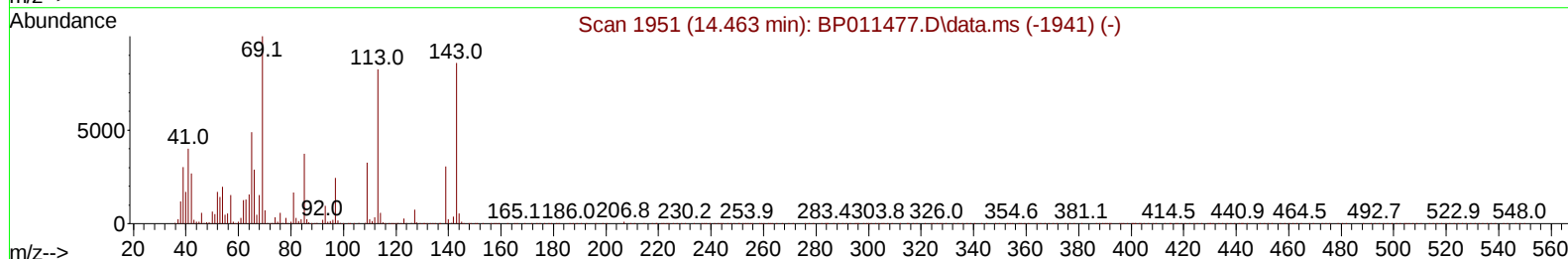
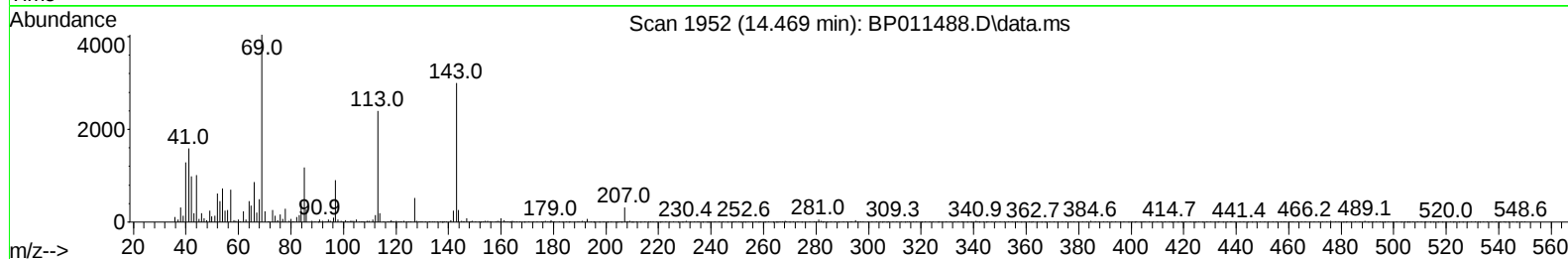
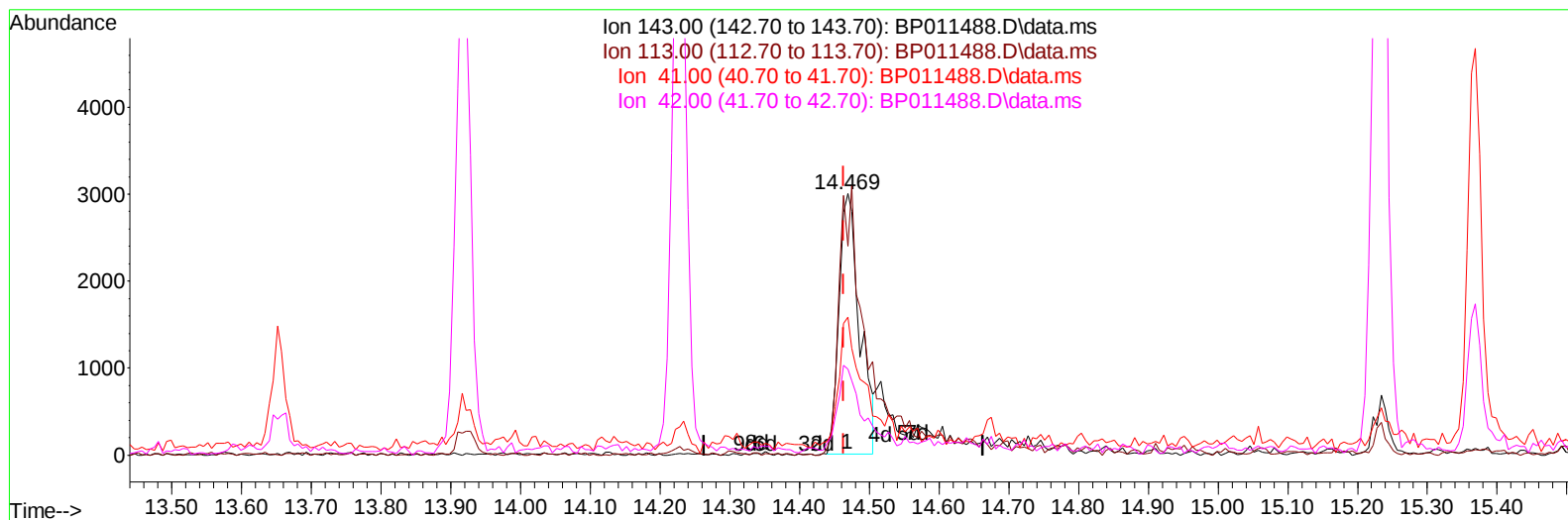
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081822\
Data File : BP011488.D
Acq On : 18 Aug 2022 18:40
Operator : CG/JU
Sample : N4221-16
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0KN8

Manual IntegrationsAPPROVED

Quant Time: Aug 18 23:18:46 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Aug 17 01:58:37 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/19/2022
Supervised By :Sohil Jodhani 08/22/2022



TIC: BP011488.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.469min (+ 0.006) 2.59 ng/u1

response 6152

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	109.60	79.86#
41.00	52.50	52.80
42.00	35.70	33.02

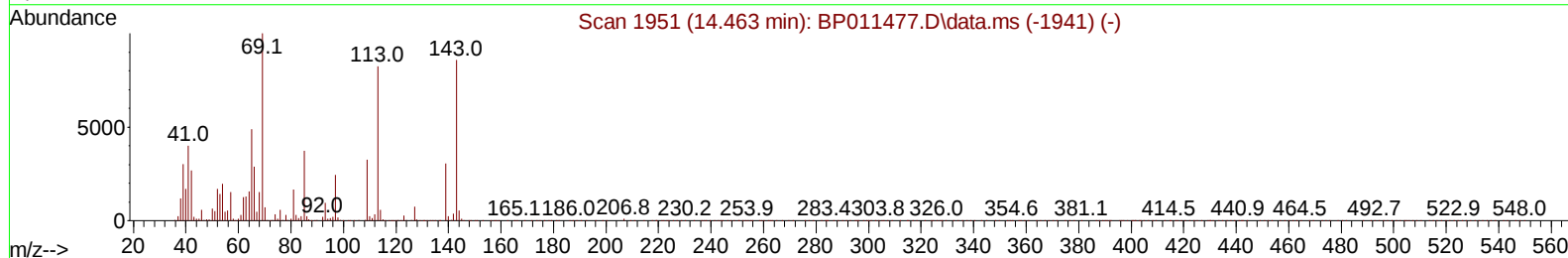
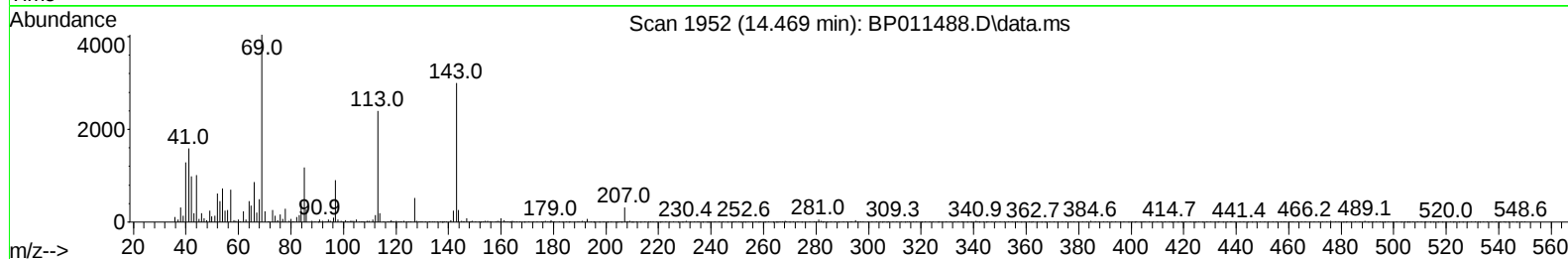
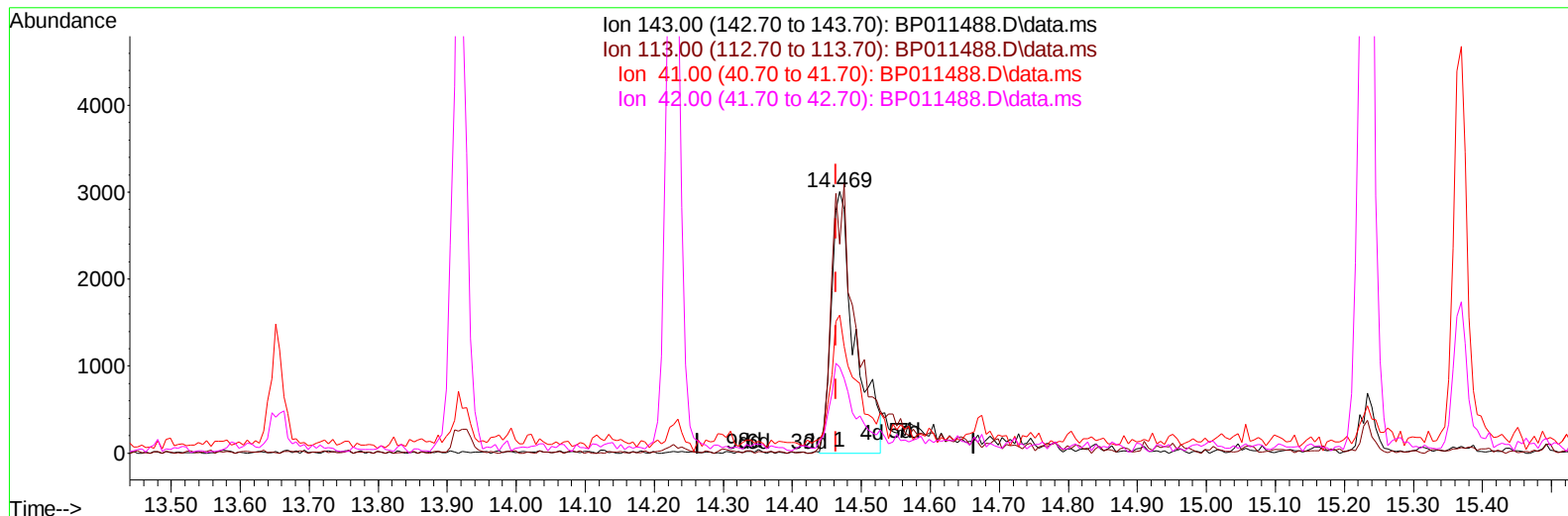
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(54) 4-Nitrophenol-d4 (S)

14.469min (+ 0.006) 3.00 ng/ul m

response 7128

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	109.60	79.86#
41.00	52.50	52.80
42.00	35.70	33.02

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.558	152	77738	20.000	ng/ul	0.00
20) Naphthalene-d8	10.346	136	319526	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.228	164	174418	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.998	188	334274	20.000	ng/ul	0.00
79) Chrysene-d12	21.127	240	303927	20.000	ng/ul	0.00
88) Perylene-d12	23.421	264	338361	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.070	96	10580	4.223	ng/uL	0.00
4) Pyridine-d5	3.487	84	52874	8.244	ng/ul	0.00
7) Phenol-d5	6.746	99	33323	4.665	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.905	67	122967	26.495	ng/ul	0.00
11) 2-Chlorophenol-d4	7.093	132	112808	21.470	ng/ul	0.00
15) 4-Methylphenol-d8	8.275	113	63495	11.105	ng/ul	0.00
21) Nitrobenzene-d5	8.722	128	69835	30.358	ng/ul	0.00
24) 2-Nitrophenol-d4	9.440	143	67886	28.997	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.969	165	104840	24.260	ng/ul	0.00
31) 4-Chloroaniline-d4	10.499	131	159742	22.206	ng/ul	0.00
46) Dimethylphthalate-d6	13.657	166	393325	31.883	ng/ul	0.00
49) Acenaphthylene-d8	13.916	160	416842	30.402	ng/ul	0.00
54) 4-Nitrophenol-d4	14.469	143	7128m	2.999	ng/ul	0.00
60) Fluorene-d10	15.234	176	320006	30.834	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.369	200	51710	28.628	ng/ul	0.00
73) Anthracene-d10	17.098	188	537934	36.336	ng/ul	0.00
81) Pyrene-d10	19.363	212	581552	37.819	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.274	264	600698	36.496	ng/ul	0.00

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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